

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110322\
 Data File : BM037346.D
 Acq On : 03 Nov 2022 14:49
 Operator : CG/JU
 Sample : PB148645BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK645

Quant Time: Nov 04 01:03:11 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM110122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 02 03:33:37 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.851	152	339413	20.000	ng/u1	0.00
20) Naphthalene-d8	10.651	136	1525740	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.486	164	1005327	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.227	188	2062515	20.000	ng/u1	0.00
79) Chrysene-d12	21.397	240	1886803	20.000	ng/u1	0.00
88) Perylene-d12	23.750	264	1574586	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	44929	5.835	ng/uL	0.00
4) Pyridine-d5	3.693	84	625637	27.220	ng/u1	0.00
7) Phenol-d5	7.022	99	821853	27.823	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.187	67	528796	29.173	ng/u1	0.00
11) 2-Chlorophenol-d4	7.381	132	681879	30.214	ng/u1	0.00
15) 4-Methylphenol-d8	8.569	113	675661	28.012	ng/u1	0.00
21) Nitrobenzene-d5	9.022	128	345753	30.337	ng/u1	0.00
24) 2-Nitrophenol-d4	9.739	143	374884	30.197	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.281	165	657556	28.729	ng/u1	0.00
31) 4-Chloroaniline-d4	10.798	131	986027	28.191	ng/u1	0.00
46) Dimethylphthalate-d6	13.904	166	2071005	29.794	ng/u1	0.00
49) Acenaphthylene-d8	14.180	160	2484091	29.726	ng/u1	0.00
54) 4-Nitrophenol-d4	14.704	143	338223	24.343	ng/u1	0.00
60) Fluorene-d10	15.480	176	1944576	31.371	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.604	200	293426	22.461	ng/u1	0.00
73) Anthracene-d10	17.327	188	3079749	32.422	ng/u1	0.00
81) Pyrene-d10	19.615	212	3381810	34.300	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.597	264	2638572	32.554	ng/u1	0.00

Target Compounds	Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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